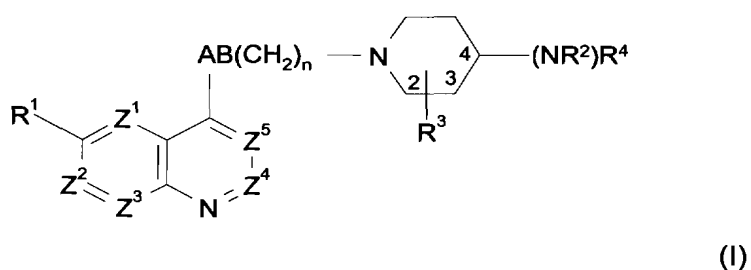


Amendments to the claims

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of claims:

1. (Currently amended) A compound of formula (I) or a pharmaceutically acceptable ~~derivative~~ salt and/or N-oxide thereof:



wherein:

one of Z¹, Z², Z³, Z⁴ and Z⁵ is N, one is CR^{1a} and the remainder are CH, CH, or one of Z¹, Z², Z³, Z⁴ and Z⁵ is CR^{1a} and the remainder are CH;

R¹ and R^{1a} are independently hydrogen; hydroxy; (C₁₋₆)alkoxy optionally substituted by (C₁₋₆)alkoxy, amino, piperidyl, guanidino or amidino any of which is optionally N-substituted by one or two (C₁₋₆)alkyl, acyl or (C₁₋₆)alkylsulphonyl groups, CONH₂, hydroxy, (C₁₋₆)alkylthio, heterocyclylthio, heterocycloxy, arylthio, aryloxy, acylthio, acyloxy or (C₁₋₆)alkylsulphonyloxy; (C₁₋₆)alkoxy-substituted(C₁₋₆)alkyl; halogen; (C₁₋₆)alkyl; (C₁₋₆)alkylthio; trifluoromethyl; trifluoromethoxy; nitro; azido; acyl; acyloxy; acylthio; (C₁₋₆)alkylsulphonyl; (C₁₋₆)alkylsulphoxide; arylsulphonyl; arylsulphoxide or an amino, piperidyl, guanidino or amidino group optionally N-substituted by one or two (C₁₋₆)alkyl, acyl or (C₁₋₆)alkylsulphonyl groups;

or when Z⁵ is CR^{1a}, R^{1a} may instead be cyano, hydroxymethyl or carboxy;

or R¹ and R^{1a} on adjacent positions may together form ethylenedioxy;

provided that when none of Z¹, Z², Z³, Z⁴ and Z⁵ is N, then R¹ is not hydrogen;

R² is hydrogen, or (C₁₋₄)alkyl or (C₂₋₄)alkenyl optionally substituted with 1 to 3 groups selected from:
amino optionally substituted by one or two (C₁₋₄)alkyl groups; carboxy; (C₁₋₄)alkoxycarbonyl; (C₁₋₄)alkylcarbonyl; (C₂₋₄)alkenyloxycarbonyl; (C₂₋₄)alkenylcarbonyl; aminocarbonyl wherein the amino group is optionally substituted by hydroxy, (C₁₋₄)alkyl, hydroxy(C₁₋₄)alkyl, aminocarbonyl(C₁₋₄)alkyl, (C₂₋₄)alkenyl, (C₁₋₄)alkylsulphonyl, trifluoromethylsulphonyl, (C₂₋₄)alkenylsulphonyl, (C₁₋₄)alkoxycarbonyl, (C₁₋₄)alkylcarbonyl, (C₂₋₄)alkenyloxycarbonyl or (C₂₋₄)alkenylcarbonyl; cyano; tetrazolyl; 2-oxo-oxazolidinyl optionally substituted by R¹⁰; 3-hydroxy-3-cyclobutene-1,2-dione-4-yl; 2,4-thiazolidinedione-5-yl; tetrazol-5-ylaminocarbonyl; 1,2,4-triazol-5-yl optionally substituted by R¹⁰; 5-oxo-1,2,4-oxadiazol-3-yl; halogen; (C₁₋₄)alkylthio; trifluoromethyl; hydroxy optionally substituted by (C₁₋₄)alkyl, (C₂₋₄)alkenyl, (C₁₋₄)alkoxycarbonyl, (C₁₋₄)alkylcarbonyl, (C₂₋₄)alkenyloxycarbonyl, or (C₂₋₄)alkenylcarbonyl; oxo; (C₁₋₄)alkylsulphonyl; (C₂₋₄)alkenylsulphonyl; ~~or~~ and (C₁₋₄)aminosulphonyl wherein the amino group is optionally substituted by (C₁₋₄)alkyl or (C₂₋₄)alkenyl;

R³ is in the 2-, 3- or 4-position and is trifluoromethyl or is in the 2-position and is oxo;
or

R³ is in the 3-position and is fluorine or amino wherein the amino group is optionally substituted by: hydroxy; (C₁₋₆)alkylsulphonyl; trifluoromethylsulphonyl; (C₂₋₆)alkenylsulphonyl; (C₁₋₆)alkylcarbonyl; (C₂₋₆)alkenylcarbonyl; (C₁₋₆)alkoxycarbonyl; (C₂₋₆)alkenyloxycarbonyl; (C₁₋₆)alkyl; or (C₂₋₆)alkenyl; wherein a (C₁₋₆)alkyl or (C₂₋₆)alkenyl moiety may be optionally substituted with up to 2 groups R¹² independently selected from:

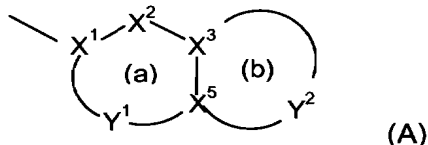
halogen; (C₁₋₆)alkylthio; trifluoromethyl; cyano; carboxy; tetrazolyl; 2-oxo-oxazolidinyl; 3-hydroxy-3-cyclobutene-1,2-dione-4-yl; 2,4-thiazolidinedione-5-yl; tetrazol-5-ylaminocarbonyl; 1,2,4-triazol-5-yl optionally substituted by R¹⁰; ~~or~~ 5-oxo-1,2,4-oxadiazol-3-yl; (C₁₋₆)alkoxycarbonyl; (C₁₋₆)alkylcarbonyl; (C₂₋₆)alkenyloxycarbonyl; (C₂₋₆)alkenylcarbonyl; hydroxy optionally substituted by (C₁₋₆)alkyl, (C₂₋₆)alkenyl, (C₁₋₆)alkylcarbonyl, (C₂₋₆)alkenylcarbonyl or aminocarbonyl wherein the amino group is optionally substituted by (C₁₋₆)alkyl, or (C₂₋₆)alkenyl; and amino optionally mono- or disubstituted by (C₁₋₆)alkoxycarbonyl, (C₁₋₆)alkylcarbonyl, (C₂₋₆)alkenyloxycarbonyl, (C₂₋₆)alkenylcarbonyl, (C₁₋₆)alkyl, (C₂₋₆)alkenyl, (C₁₋₆)alkylsulphonyl, (C₂₋₆)alkenylsulphonyl or aminocarbonyl wherein the amino group is optionally substituted by (C₁₋₆)alkyl or (C₂₋₆)alkenyl;

in addition when R³ is disubstituted with a hydroxy or amino containing substituent and a carboxy containing substituent these may together form a cyclic ester or amide linkage, respectively;

R⁴ is a group -U-R⁵ where

U is selected from CO, SO₂ and CH₂ and

R⁵ is an optionally substituted bicyclic carbocyclic or heterocyclic ring system (A):



containing up to four heteroatoms in each ring in which

ring (a) is aromatic and ring (b) is non-aromatic;

X¹ is C or N;

X² is N, NR¹³, O, S(O)_x, CO or CR¹⁴;

X³ and X⁵ are independently N or C;

Y¹ is a 0 to 4 atom linker group each atom of which is independently selected from N, NR¹³, O, S(O)_x, CO and CR¹⁴;

Y² is a 2 to 6 atom linker group, each atom of Y² being independently selected from N, NR¹³, O, S(O)_x, CO, CR¹⁴ and CR¹⁴R¹⁵;

each of R¹⁴ and R¹⁵ is independently selected from: H; (C₁₋₄)alkylthio; halo; carboxy(C₁₋₄)alkyl; halo(C₁₋₄)alkoxy; halo(C₁₋₄)alkyl; (C₁₋₄)alkyl; (C₂₋₄)alkenyl; (C₁₋₄)alkoxycarbonyl; formyl; (C₁₋₄)alkylcarbonyl; (C₂₋₄)alkenyloxycarbonyl; (C₂₋₄)alkenylcarbonyl; (C₁₋₄)alkylcarbonyloxy; (C₁₋₄)alkoxycarbonyl(C₁₋₄)alkyl; hydroxy; hydroxy(C₁₋₄)alkyl; mercapto(C₁₋₄)alkyl; (C₁₋₄)alkoxy; trifluoromethoxy; nitro; cyano; carboxy; amino or aminocarbonyl optionally substituted as for corresponding substituents in R³; (C₁₋₄)alkylsulphonyl; (C₂₋₄)alkenylsulphonyl; α -aminosulphonyl wherein the amino group is optionally mono- or di-substituted by (C₁₋₄)alkyl or (C₂₋₄)alkenyl; aryl; aryl(C₁₋₄)alkyl; and aryl(C₁₋₄)alkoxy;

each R¹³ is independently H; trifluoromethyl; (C₁₋₄)alkyl optionally substituted by hydroxy, (C₁₋₆)alkoxy, (C₁₋₆)alkylthio, halo or trifluoromethyl; (C₂₋₄)alkenyl; aryl; aryl (C₁₋₄)alkyl; arylcarbonyl; heteroarylcarbonyl; (C₁₋₄)alkoxycarbonyl; (C₁₋₄)alkylcarbonyl; formyl; (C₁₋₆)alkylsulphonyl; or aminocarbonyl wherein the amino group is optionally substituted by (C₁₋₄)alkoxycarbonyl, (C₁₋₄)alkylcarbonyl, (C₂₋₄)alkenyloxycarbonyl, (C₂₋₄)alkenylcarbonyl, (C₁₋₄)alkyl or (C₂₋₄)alkenyl and optionally further substituted by (C₁₋₄)alkyl or (C₂₋₄)alkenyl;

each x is independently 0, 1 or 2

n is 0 and AB is NR^{11}CO , $\text{CO-CR}^8\text{R}^9$, $\text{CR}^6\text{R}^7\text{-CO}$, $\text{NHR}^{11}\text{SO}_2$, $\text{CR}^6\text{R}^7\text{-SO}_2$ or $\text{CR}^6\text{R}^7\text{-CR}^8\text{R}^9$, provided that R^8 and R^9 are not optionally substituted hydroxy or amino and R^6 and R^8 do not represent a bond;
or n is 1 and AB is NR^{11}CO , $\text{CO-CR}^8\text{R}^9$, $\text{CR}^6\text{R}^7\text{-CO}$, $\text{NR}^{11}\text{SO}_2$, CONR^{11} , $\text{CR}^6\text{R}^7\text{-CR}^8\text{R}^9$, $\text{O-CR}^8\text{R}^9$ or $\text{NR}^{11}\text{-CR}^8\text{R}^9$;
each of R^6 , R^7 , R^8 and R^9 is independently selected from: hydrogen; (C_{1-6}) alkoxy; (C_{1-6}) alkylthio; halo; trifluoromethyl; azido; (C_{1-6}) alkyl; (C_{2-6}) alkenyl; (C_{1-6}) alkoxycarbonyl; (C_{1-6}) alkylcarbonyl; (C_{2-6}) alkenyloxycarbonyl; (C_{2-6}) alkenylcarbonyl; hydroxy, amino or aminocarbonyl optionally substituted as for corresponding substituents in R^3 ; (C_{1-6}) alkylsulphonyl; (C_{2-6}) alkenylsulphonyl; or and aminosulphonyl wherein the amino group is optionally substituted by (C_{1-6}) alkyl or (C_{2-6}) alkenyl;
or when $n=1$ R^6 and R^8 together represent a bond and R^7 and R^9 are as above defined;
or R^6 and R^7 or R^8 and R^9 together represent oxo;

R^{10} is selected from (C_{1-4}) alkyl; (C_{2-4}) alkenyl and aryl any of which may be optionally substituted by a group R^{12} as defined above; carboxy; aminocarbonyl wherein the amino group is optionally substituted by hydroxy, (C_{1-6}) alkyl, (C_{2-6}) alkenyl, (C_{1-6}) alkylsulphonyl, trifluoromethylsulphonyl, (C_{2-6}) alkenylsulphonyl, (C_{1-6}) alkoxycarbonyl, (C_{1-6}) alkylcarbonyl, (C_{2-6}) alkenyloxycarbonyl or (C_{2-6}) alkenylcarbonyl and optionally further substituted by (C_{1-6}) alkyl or (C_{2-6}) alkenyl; and

R^{11} is hydrogen; trifluoromethyl, (C_{1-6}) alkyl; (C_{2-6}) alkenyl; (C_{1-6}) alkoxycarbonyl; (C_{1-6}) alkylcarbonyl; or aminocarbonyl wherein the amino group is optionally substituted by (C_{1-6}) alkoxycarbonyl, (C_{1-6}) alkylcarbonyl, (C_{2-6}) alkenyloxycarbonyl, (C_{2-6}) alkenylcarbonyl, (C_{1-6}) alkyl or (C_{2-6}) alkenyl and optionally further substituted by (C_{1-6}) alkyl or (C_{2-6}) alkenyl;

or where one of R^3 and R^6 , R^7 , R^8 or R^9 contains a carboxy group and the other contains a hydroxy or amino group they may together form a cyclic ester or amide linkage.

2. (Original) A compound according to claim 1 wherein Z^5 is CH, C-Cl or N, Z^3 is CH or CF and Z^1 , Z^2 and Z^4 are each CH, or Z^1 is N, Z^3 is CH and Z^2 and Z^4 are each CH and Z^5 is CH or C-Cl.

3. (Previously presented) A compound according to claim 1 wherein R^1 is methoxy and R^{1a} is H or when Z^3 is CR^{1a} it may be C-F or when Z^5 is CR^{1a} it may be C-F or C-Cl.
4. (Previously presented) A compound according to claim 1 wherein R^2 is hydrogen, carboxymethyl, hydroxyethyl, aminocarbonylmethyl, ethoxycarbonylmethyl, ethoxycarbonylallyl or carboxyallyl.
5. (Previously presented) A compound according to claim 1 wherein R^3 is CF_3 , fluoro, oxo or amino unsubstituted or substituted by (C_{1-6}) alkyl or (C_{2-6}) alkenyl.
6. (Previously presented) A compound according to claim 1 wherein n is 0 and either A is CH_2 or $CHOH$ and B is CH_2 or A is NH and B is CO .
7. (Previously presented) A compound according to claim 1 wherein $-U-$ is $-CH_2-$.
8. (Currently Amended) A compound according to claim 1 wherein in the heterocyclic ring (A) ring (a) is selected from optionally substituted benzo and pyrido and Y^2 has 3-5 atoms including a heteroatom bonded to X^5 selected from NR^{13} , O or and S, where R^{13} is other than hydrogen, and $NHCO$ bonded via N to X^3 , or O or NH bonded to X^3 .
9. (Currently Amended) A compound according to claim 1 wherein R^5 is selected from:
 - 4H-benzo[1,4] oxazin-3-one-6-yl;
 - 4H-benzo[1,4] thiazin-3-one-6-yl;
 - 2,3-dihydro-[1,4]dioxino[2,3-c]pyridin-7-yl;
 - 3-oxo-3,4-dihydro-2H-pyrido[3,2-b][1,4]thiazin-6-yl;
 - 3-oxo-3,4-dihydro-2H-pyrido[3,2-b][1,4]oxazin-6-yl;
 - 7-chloro-3-oxo-3,4-dihydro-2H-pyrido[3,2-b][1,4]oxazin-6-yl; and
 - 7-chloro-3-oxo-3,4-dihydro-2H-pyrido[3,2-b][1,4]thiazin-6-yl.
10. (Currently amended) A compound according to claim 1 selected from:
6-({2S,4S)-1-[(R)-2-Hydroxy-2-(6-methoxy-[1,5]naphthyridin-4-yl)ethyl]-2-(trifluoromethyl)piperidin-4-ylamino)methyl)-4H-benzo[1,4]thiazin-3-one;

6-((3*R*,4*S*)-1-[(*R*)-2-Hydroxy-2-(6-methoxy-[1,5]naphthyridin-4-yl)ethyl]-3-(trifluoromethyl)piperidin-4-ylamino)methyl)-4*H*-benzo[1,4]thiazin-3-one;
6-((1-[(*R*)-2-Hydroxy-2-(6-methoxy-[1,5]naphthyridin-4-yl)ethyl]-4-(trifluoromethyl)piperidin-4-ylamino)methyl)-4*H*-benzo[1,4]thiazin-3-one;
6-((1-[(*R*)-2-Hydroxy-2-(6-methoxyquinolin-4-yl)ethyl]-2-oxopiperidin-4-ylamino)methyl)-4*H*-benzo[1,4]thiazin-3-one;
6-[(((3*S*,4*R*)-3-Fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-benzo[1,4]thiazin-3-one and 6-[(((3*R*,4*S*)-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-benzo[1,4]thiazin-3-one ;
6-((cis-3-Fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one Diastereoisomer 1 ;
6-((cis-3-Fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]-piperidin-4-ylamino)-methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one Diastereoisomer 2 ;
7-Chloro-6-((cis-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one Diastereoisomer 1 ;
7-Chloro-6-((cis-3-Fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one Diastereoisomer 2 ;
6-((cis-3-Fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one Diastereoisomer 1 ;
6-((cis-3-Fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one Diastereoisomer 2 ;
7-Chloro-6-[(((3*S*,4*R*)-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one and 7-chloro-6-[(((3*R*,4*S*)-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one ;
7-Fluoro-6-(((3*S*,4*R*)-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one and 7-fluoro-6-[(((3*R*,4*S*)-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one;
7-(((3*S*,4*R*)-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-1*H*-pyrido[2,3-*b*][1,4]thiazin-2-one and 7-(((3*R*,4*S*)-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-1*H*-pyrido[2,3-*b*][1,4]thiazin-2-one;
7-Chloro-6-[(((3*S*,4*R*)-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one and 7-chloro-6-[(((3*R*,4*S*)-3-fluoro-1-[(*R*)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one;

6-(((3S,4S)-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one and 6-(((3*R*,4*R*)-3-fluoro-1-[(R)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one;

6-(((3S,4S)-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one and 6-(((3*R*,4*R*)-3-fluoro-1-[(R)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one;

7-Fluoro-6-(((3S,4S)-3-fluoro-1-[(R)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one and 7-Fluoro-6-(((3*R*,4*R*)-3-fluoro-1-[(R)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one;

6-(((3S,4S)-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-1*H*-pyrido[2,3-*b*][1,4]thiazin-3-one and 6-(((3*R*,4*R*)-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-1*H*-pyrido[2,3-*b*][1,4]thiazin-3-one;

6-((cis-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one Diastereoisomer 1;

6-((cis-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one Diastereoisomer 2;

6-((cis-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one Diastereoisomer 1;

6-((cis-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one Diastereoisomer 2;

7-Chloro-6-((cis-3-fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one Diastereoisomer 1;

7-Chloro-6-((cis-3-fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one Diastereoisomer 2;

6-((cis-3-Fluoro-1-[(R)-2-hydroxy-2-(8-fluoro-6-methoxy-quinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]oxazin-3-one Diastereoisomer 1;

6-((cis-3-Fluoro-1-[(R)-2-hydroxy-2-(8-fluoro-6-methoxy-quinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one Diastereoisomer 1;

6-(((3*R*,4*S*)-1-[2-(3-Chloro-6-methoxy-quinolin-4-yl)-ethyl]-3-fluoro-piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one and 6-(((3*S*,4*R*)-1-[2-(3-Chloro-6-methoxy-quinolin-4-yl)-ethyl]-3-fluoro-piperidin-4-ylamino)methyl)-4*H*-pyrido[3,2-*b*][1,4]thiazin-3-one;

6-(((3R,4S)-3-Fluoro-1-[2-(6-methoxy-[1,5]naphthyridin-4-yl)-ethyl]-piperidin-4-ylamino)-methyl)-4H-pyrido[3,2-b][1,4]thiazin-3-one and 6-(((3S,4R)-3-fluoro-1-[2-(6-methoxy-[1,5]naphthyridin-4-yl)-ethyl]-piperidin-4-ylamino)-methyl)-4H-pyrido[3,2-b][1,4]thiazin-3-one;

6-(((3S,4R)-3-Fluoro-1-[(S)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4H-pyrido[3,2-b][1,4]thiazin-3-one and 6-(((3R,4S)-3-fluoro-1-[(S)-2-hydroxy-2-(6-methoxyquinolin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4H-pyrido[3,2-b][1,4]thiazin-3-one;

6-(((3R,4S)-1-[2-(2,3-Dihydro-[1,4]dioxino[2,3-f]quinolin-10-yl)-ethyl]-3-fluoro-piperidin-4-ylamino)-methyl)-4H-pyrido[3,2-b][1,4]thiazin-3-one and 6-(((3S,4R)-1-[2-(2,3-dihydro-[1,4]dioxino[2,3-f]quinolin-10-yl)-ethyl]-3-fluoro-piperidin-4-ylamino)-methyl)-4H-pyrido[3,2-b][1,4]thiazin-3-one;

6-(((3R,4S)-1-[2-(6,8-Difluoro-quinolin-4-yl)-ethyl]-3-fluoro-piperidin-4-ylamino)-methyl)-4H-pyrido[3,2-b][1,4]thiazin-3-one and 6-(((3S,4R)-1-[2-(6,8-difluoro-quinolin-4-yl)-ethyl]-3-fluoro-piperidin-4-ylamino)-methyl)-4H-pyrido[3,2-b][1,4]thiazin-3-one;

6-(((3S,4R)-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4H-benzo[1,4]thiazin-3-one and 6-(((3R,4S)-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4H-benzo[1,4]thiazin-3-one;

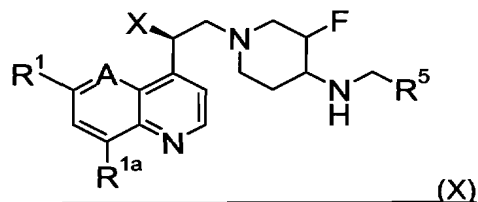
6-(((cis-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4H-benzo[1,4]thiazin-3-one Faster running Diastereoisomer;

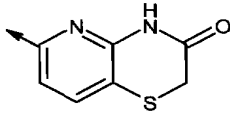
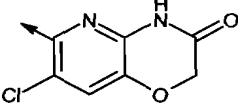
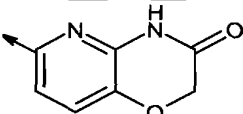
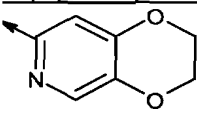
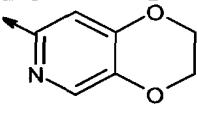
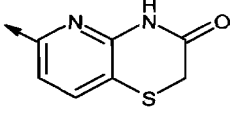
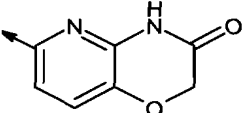
6-(((cis-3-Fluoro-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin-4-yl)-ethyl]piperidin-4-ylamino)methyl)-4H-benzo[1,4]thiazin-3-one Slower-running Diastereoisomer;

6-((2S,4S)-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin-4-yl)ethyl]-2-(trifluoromethyl)piperidin-4-ylamino)methyl)-4H-pyrido[1,4]thiazin-3-one ;

6-((2S,4R)-1-[(R)-2-hydroxy-2-(6-methoxy-[1,5]naphthyridin-4-yl)ethyl]-2-(trifluoromethyl)piperidin-4-ylamino)methyl)-4H-pyrido[1,4]thiazin-3-one;

and the following tabulated compounds of formula (X):



<u>Isomeric form</u>	<u>A</u>	<u>R¹</u>	<u>R^{1a}</u>	<u>X</u>	<u>R⁵</u>
<u>Enantiomer 2</u>	<u>CH</u>	<u>MeO</u>	<u>F</u>	<u>OH</u>	<u>6-[4H-pyrido[3,2-b][1,4]thiazin-3-one]</u> 
<u>Enantiomer 1</u>	<u>CH</u>	<u>MeO</u>	<u>F</u>	<u>OH</u>	<u>6-[7-chloro-4H-pyrido[3,2-b][1,4]oxazin-3-one]</u> 
<u>Enantiomer 2</u>	<u>CH</u>	<u>MeO</u>	<u>F</u>	<u>OH</u>	<u>6-[4H-pyrido[3,2-b][1,4]oxazin-3-one]</u> 
<u>Enantiomer 2</u>	<u>CH</u>	<u>MeO</u>	<u>H</u>	<u>OH</u>	<u>7-[2,3-dihydro-[1,4]dioxino[2,3-c]pyridine]</u> 
<u>Enantiomer 1</u>	<u>CH</u>	<u>MeO</u>	<u>H</u>	<u>OH</u>	<u>7-[2,3-dihydro-[1,4]dioxino[2,3-c]pyridine]</u> 
<u>Enantiomer 2</u>	<u>N</u>	<u>MeO</u>	<u>H</u>	<u>H</u>	<u>6-[4H-pyrido[3,2-b][1,4]thiazin-3-one]</u> 
<u>Racemic</u>	<u>CH</u>	<u>F</u>	<u>F</u>	<u>H</u>	<u>6-[4H-pyrido[3,2-b][1,4]oxazin-3-one]</u> 

or a pharmaceutically acceptable ~~derivative~~ salt and/or N-oxide thereof.

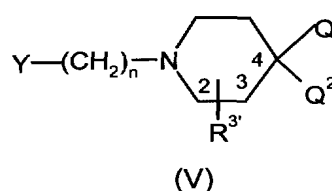
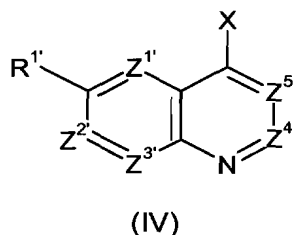
11. (Currently amended) A method of treatment of bacterial infections in mammals, ~~particularly in man~~, which method comprises the administration to a mammal in need of such treatment an effective amount of a compound according to claim 1.

Claims 12 and 13 (Cancelled).

14. (Original) A pharmaceutical composition comprising a compound according to claim 1, and a pharmaceutically acceptable carrier.

Claims 15 and 16 (Cancelled).

17. (Currently amended) A process for preparing a compound of formula (I) according to claim 1, or a pharmaceutically acceptable ~~derivative salt and/or N-oxide~~ thereof, which process comprises reacting a compound of formula (IV) with a compound of formula (V):

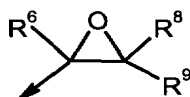


wherein n is as defined in formula (I); $Z^{1'}$, $Z^{2'}$, $Z^{3'}$, $Z^{4'}$, $Z^{5'}$, $R^{1'}$, and $R^{3'}$ are Z^1 , Z^2 , Z^3 , Z^4 , Z^5 , R^1 , and R^3 are as defined in formula (I) or groups convertible thereto; Q^1 is $NR^{2'}R^{4'}$ or a group convertible thereto wherein $R^{2'}$ and $R^{4'}$ are R^2 and R^4 as defined in formula (I) or groups convertible thereto and Q^2 is H or $R^{3'}$ or Q^1 and Q^2 together form an optionally protected oxo group;

- (i) X is A'-COW, Y is H and n is 0;
- (ii) X is $CR^6=CR^8R^9$, Y is H and n is 0;
- (iii) X is oxirane, Y is H and n is 0;
- (iv) X is $N=C=O$ and Y is H and n is 0;
- (v) one of X and Y is CO_2R^Y and the other is $CH_2CO_2R^X$;
- (vi) X is CHR^6R^7 and Y is $C(=O)R^9$;
- (vii) X is $CR^7=PR^{Z_3}$ and Y is $C(=O)R^9$ and $n=1$;
- (viii) X is $C(=O)R^7$ and Y is $CR^9=PR^{Z_3}$ and $n=1$;
- (ix) Y is COW and X is $NHR^{11'}$ or $NR^{11'}COW$ and $n=0$ or 1 or when $n=1$ X is COW and Y is $NHR^{11'}$ or $NR^{11'}COW$;
- (x) X is $NHR^{11'}$ and Y is $C(=O)R^8$ and $n=1$;
- (xi) X is $NHR^{11'}$ and Y is CR^8R^9W and $n=1$;
- (xii) X is $NR^{11'}COCH_2W$ or $NR^{11'}SO_2CH_2W$ and Y is H and $n=0$;
- (xiii) X is $CR^6R^7SO_2W$ and Y is H and $n=0$;
- (xiv) X is W or OH and Y is CH_2OH and n is 1;
- (xv) X is $NHR^{11'}$ and Y is SO_2W or X is $NR^{11'}SO_2W$ and Y is H, and n is 0;

(xvi) X is W and Y is CONHR^{11'};

in which W is a leaving group, ~~e.g. halo or imidazolyl~~; R^X and R^Y are (C₁₋₆)alkyl; R^Z is aryl or (C₁₋₆)alkyl; A' and NR^{11'} are A and NR¹¹ as defined in formula (I), or groups convertible thereto; and oxirane is:



wherein R⁶, R⁸ and R⁹ are as defined in formula (I);

and thereafter optionally or as necessary converting Q¹ and Q² to NR^{2'}R^{4'};

converting A', Z^{1'}, Z^{2'}, Z^{3'}, Z^{4'}, Z^{5'}, R^{1'}, R^{2'}, R^{3'}, R^{4'} and ~~NR^{11'}~~ to NR^{11'} to A, Z¹,

Z², Z³, Z⁴, Z⁵, R¹, R², R³, R⁴ and NR¹¹; converting A-B to other A-B,

interconverting R¹, R², R³ and/or R⁴, and/or forming a pharmaceutically acceptable

~~derivative~~ salt and/or N-oxide thereof.

18. (New) A compound according to claim 1 wherein R³ is fluoro.